

Data Path : Z:\svoasrv\HPCHEM1\BNA\_M\Data\BM042621\  
 Data File : BM029661.D  
 Acq On : 26 Apr 2021 22:02  
 Operator : CG/JU  
 Sample : M2170-12  
 Misc :  
 ALS Vial : 14 Sample Multiplier: 1

Instrument :  
 BNA\_M  
 ClientSampleId :  
 CS-1

Manual Integrations  
 APPROVED

mohammad  
 4/27/2021 12:11:28 PM

Quant Time: Apr 27 01:33:54 2021  
 Quant Method : Z:\svoasrv\HPCHEM1\BNA\_M\Methods\8270-BM042121.M  
 Quant Title : ASP BNA STANDARDS FOR 5 POINT CALIBRATION  
 QLast Update : Mon Apr 26 15:01:01 2021  
 Response via : Initial Calibration

| Compound                    | R.T.   | QIon | Response | Conc   | Units | Dev(Min)  |
|-----------------------------|--------|------|----------|--------|-------|-----------|
| -----                       |        |      |          |        |       |           |
| Internal Standards          |        |      |          |        |       |           |
| 1) 1,4-Dichlorobenzene-d4   | 7.616  | 152  | 60472    | 20.00  | ng    | 0.00      |
| 21) Naphthalene-d8          | 10.398 | 136  | 213225   | 20.00  | ng    | 0.00      |
| 39) Acenaphthene-d10        | 14.263 | 164  | 142065   | 20.00  | ng    | 0.00      |
| 64) Phenanthrene-d10        | 17.009 | 188  | 320467   | 20.00  | ng    | 0.00      |
| 76) Chrysene-d12            | 21.192 | 240  | 371630   | 20.00  | ng    | 0.00      |
| 86) Perylene-d12            | 23.380 | 264  | 398114   | 20.00  | ng    | 0.00      |
| System Monitoring Compounds |        |      |          |        |       |           |
| 5) 2-Fluorophenol           | 5.222  | 112  | 460084   | 152.53 | ng    | 0.00      |
| 7) Phenol-d6                | 6.798  | 99   | 510370   | 113.28 | ng    | 0.00      |
| 23) Nitrobenzene-d5         | 8.775  | 82   | 432096   | 104.37 | ng    | 0.00      |
| 42) 2,4,6-Tribromophenol    | 15.762 | 330  | 346835   | 176.53 | ng    | 0.00      |
| 45) 2-Fluorobiphenyl        | 12.880 | 172  | 1160264  | 110.85 | ng    | 0.00      |
| 79) Terphenyl-d14           | 19.645 | 244  | 2113226  | 107.81 | ng    | 0.00      |
| Target Compounds            |        |      |          |        |       |           |
| 10) Phenol                  | 6.828  | 94   | 14579    | 3.34   | ng    | Qvalue 88 |
| 32) Benzoic acid            | 9.710  | 122  | 1679m    | 7.80   | ng    |           |
| 50) Dimethylphthalate       | 13.727 | 163  | 58399    | 5.44   | ng    | 98        |
| -----                       |        |      |          |        |       |           |

(#) = qualifier out of range (m) = manual integration (+) = signals summed

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